## Amendments to the Claims

Please amend the claims as follows (the changes in these claims are shown with strikethrough for deleted text and <u>underlines</u> for added text). A complete listing of the claims is listed below with proper claim identifiers. This listing of claims will replace all prior versions, and listings, of claims in the application.

## What is claimed is:

1. (Currently Amended) A modulator of the structure (I), or a salt thereof:

$$\mathbb{R}^{4} \xrightarrow{\mathbb{R}^{3}} \mathbb{Q}^{\mathbb{Z}} \mathbb{Q}^{\mathbb$$

where m is an integer from 1 to 5;

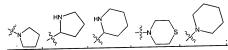
each Y is independently selected from the group consisting of hydrogen, halogen, -CN, -NO<sub>2</sub>, -OH, -OR', -C(O)R', -CO<sub>2</sub>R', -O(CO)R', -C(O)NR'R', -OC(O)NR'R', -SR', -SOR, -SO<sub>2</sub>R', -SO<sub>2</sub>NR'R', -NR'R'', -NR'C(O)R', -NR'C(O)<sub>2</sub>R', -NR'SO<sub>2</sub>R', -NR'(CO)NR'R', unsubstituted or substituted C<sub>1-8</sub> alkyl, unsubstituted or substituted C<sub>2-8</sub> alkenyl, unsubstituted or substituted 3- to 10-membered heteroaryl, and unsubstituted or substituted 3- to 10-membered heterocyclyl;

where each R', R" and R" are independently hydrogen, halogen, unsubstituted or substituted  $C_{1-8}$  alkyl, unsubstituted or substituted  $C_{6-10}$  aryl, unsubstituted or substituted 5- to 10-membered heteroaryl, and unsubstituted or substituted 3- to 10-membered heterocyclyl;

n is 0, 1, 2 or 3;

R<sup>1</sup>-and-R<sup>2</sup> are each independently substituted or unsubstituted alkyl or hydrogen, or Z in combination with R<sup>1</sup> and R<sup>2</sup> form a substituted or unsubstituted 5- to 8-membered ring comprising at least one nitrogen and 0 to 3 additional heteroatoms;

## Z is a substituted or unsubstituted group of the formulae:



R6 is alkyl, hydrogen, or halogen; and

 $R^3,\,R^4,\,$  and  $R^5$  are each independently selected from the group consisting of hydrogen, halogen, -CN, -NO<sub>2</sub>, -OH, -OR', -C(O)R', -CO<sub>2</sub>R', -O(CO)R', -C(O)NR'R", -OC(O)NR'R", -SR', -SOR, -SO<sub>2</sub>R', -SO<sub>2</sub>NR'R", -NR'R", -NR'C(O)R", -NR'C(O)<sub>2</sub>R", -NR'SO<sub>2</sub>R', -NR'C(O)R'R", -NR'C(O)R", -NR'C(O)R'R", -NR'SO<sub>2</sub>R', -NR'C(O)R'R", unsubstituted or substituted  $C_{2:8}$  alkenyl, unsubstituted or substituted  $C_{2:8}$  alkenyl, unsubstituted or substituted or substituted

- 2. (Previously presented) The modulator of claim 1, where  $\mathsf{R}^6$  is hydrogen.
- (Currently amended) The modulator of claim 1, where R<sup>6</sup> is substituted or unsubstituted C<sub>1-8</sub> alkyl.
- (Previously presented) The modulator of claim 1, where R<sup>6</sup> is halogen.
- (Currently amended) The modulator of claim 1, where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup>
  are each independently selected from the group consisting of hydrogen, OR', and substituted or unsubstituted C<sub>1.8</sub> alkyl.
- (Currently amended) The modulator of claim 1, where R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are each independently selected from the group consisting of –OR' and hydrogen.
- 7. (Previously presented) The modulator of claim 1, where  $R^3$ ,  $R^4$ , and  $R^5$  are each –OR', where R' is substituted  $C_{1:8}$  alkyl.
- (Previously presented) The modulator of claim 1, where R<sup>4</sup> and R<sup>5</sup> together with the atom which they substitute form substituted or unsubstituted 5- to 6-membered heterocyclyl containing 1 to 2 oxygen atoms.
- 9. (Previously presented) The modulator of claim 1, where Z is  $CHR^1R^2$  and where  $R^1$  and  $R^2$  together with Z form  $C_{3-10}$  cycloalkyl with 0 to 3 substituents selected from the group consisting of halogen, -CN, -NO<sub>2</sub>, -OH, -OR', -C(O)R', -CO<sub>2</sub>R', -O(CO)R', -C(O)NR'R'', -OC(O)NR'R'', -SR', -SOR', -SO<sub>2</sub>R', -SO<sub>2</sub>NR'R'', -NR'R'', -NR'C(O)R'', -NR'C(O)<sub>2</sub>R', -NR'SO<sub>2</sub>R', -NR'C(O)NR'R'', unsubstituted or substituted  $C_{1-8}$  alkyl, unsubstituted or substituted  $C_{2-8}$  alkenyl, unsubstituted  $C_{3-8}$  cycloalkyl, unsubstituted or substituted  $C_{8-10}$  aryl, unsubstituted or substituted  $C_{3-8}$  cycloalkyl, unsubstituted or substituted  $C_{3-10}$  aryl, unsubstituted  $C_{3-10}$  aryl,
- 10. (Canceled)

- 11. (Canceled)
- 12. (Canceled)
- 13. (Canceled)
- 14. (Previously presented) The modulator of claim 1, where Z is a substituted or unsubstituted group of the formula:

15. (Previously presented) The modulator of claim 1, where Z is a substituted or unsubstituted group of the formula:

16. (Previously presented) The modulator of claim 1, where Z is a substituted or unsubstituted group of the formula:

17. (Previously presented) The modulator of claim 1, where Z is a substituted or unsubstituted group of the formula:

18. (Previously presented) The modulator of claim 1, where Z is a substituted or unsubstituted group of the formula:

19. (Previously presented) The modulator of claim 1, where Z is a substituted or unsubstituted group of the formula:

where  $R^7$  is selected from the group consisting of hydrogen, -C(O)R',  $-CO_2R'$ , -C(O)NR'R'',  $-SO_2R'$ , unsubstituted or substituted  $C_{1-10}$  alkyl, unsubstituted or substituted  $C_{2-10}$  alkenyl, unsubstituted or substituted or substituted or substituted or substituted or substituted or substituted  $C_{2-10}$  alkynyl, unsubstituted or substituted  $C_{3-10}$  cycloalkyl, unsubstituted or substituted  $C_{6-10}$  aryl, C6-10 aryloxy unsubstituted or substituted 5- to  $C_{6-10}$  aryl, and unsubstituted or substituted 3- to  $C_{6-10}$  aryloxy unsubstituted or substituted  $C_{6-10}$  aryloxy unsubstituted  $C_{6-10}$  aryloxy unsubstituted

- 20. (Previously presented) The modulator of claim 1, where  $R^7$  is substituted or unsubstituted  $C_{1-10}$  alkyl, substituted or unsubstituted  $C_{1-10}$  alkoxy, substituted or unsubstituted aryloxy, or substituted or unsubstituted  $C_{3-10}$  cycloalkyl.
- 21. (Original) The modulator of claim 1, where n is 1, 2, or 3.
- 22. (Currently amended) The modulator of claim 1, where m is 1 or 2, and each Y is a halogen.
- 23. (Previously presented) The modulator of claim 1, where m is 0.
- 24. (Currently amended) The modulator of claim 1, where substituted alkyl, substituted alkyl, substituted alkynyl and substituted cycloalkyl can each independently be substituted 1 to 3 times with halogen, -OR', -NR'R'', -SR', -SIR'R''R''', -OC(O)R', -C(O)R', -CO\_2R', -CONR'R'', -OC(O)NR'R'', -NR'C(O)R', -NR'C(O)R'', -NR'C(O)\_2R', -S(O)\_2R', -S(O)\_2R', -S(O)\_2R'', -S(O)\_2R''

unsubstituted aryl, unsubstituted heteroaryl, unsubstituted or substituted heterocyclyl.

- 25. (Currently amended) The modulator of claim 1, where substituted aryl and substituted heteroaryl can each independently be substituted 1 to 3 times with halogen, unsubstituted or substituted alkyl, unsubstituted or substituted alkenyl, unsubstituted or substituted districted alkenyl, unsubstituted or substituted cycloalkyl, -OR', oxo (=0 or -O), -OC(O)R', -NR'R", -SR', -R', -CN, -NO<sub>2</sub>, -CO<sub>2</sub>R', -CONR'R", -C(O)R', -OC(O)NR'R", -NR'C(O)R', -NR'C(O)R', -NR'C(O)R', -NR'C(O)R', -NR'C(O)R', -NR'C(O)R', -NR'C(O)R', -NR'C(O)R', -S(O)R', -S(O)R', -S(O)R', -NR'C(O)R', -NR'C(O)R', alkyl, unsubstituted C3-6 cycloalkyl, unsubstituted C2-8 alkynl, unsubstituted heteroaryl, unsubstituted heterocyclyl.
- 26. (Currently amended) The modulator of claim 1, where substituted heterocyclyl can be substituted 1 to 3 times with halogen, unsubstituted or substituted alkyl, unsubstituted or substituted alkenyl, unsubstituted or substituted cycloalkyl, -OR', oxo (=O or -O), -OC(O)R', -NR'R", -SR', -R', -CN, -NO<sub>2</sub>, -OC(O)NR'R", -NR"C(O)R', -NR"C(O)R', -NR"-C(O)NR"R", -NH-C(NH<sub>2</sub>)=NH, -NR'C(O)R", and -N<sub>3</sub>, where R', R" and R" each independently hydrogen, halogen, unsubstituted C<sub>1-8</sub> alkyl, unsubstituted or C<sub>3-6</sub> cycloalkyl, unsubstituted C<sub>2-8</sub> alkenyl, unsubstituted heterocyclyl.
- 27. (Currently amended) A modulator having the structure (II):

$$\mathbb{R}^3$$
  $\mathbb{N}\mathbb{R}^7$   $\mathbb{N}^7$ 

where n=0-4

where each Y is independently hydrogen or halogen;

R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are each independently R<sup>3</sup>, R<sup>4</sup>, and R<sup>5</sup> are each independently selected from the group consisting of hydrogen, halogen, and -OR';

or any two of  $R^3,\,R^4,$  and  $R^5,$  together with the atoms which they substituted, form unsubstituted or substituted 3- to 10-membered heterocyclyl; and

 $R^7$  is selected from the group consisting of hydrogen, -C(O)R', -CO<sub>2</sub>R', -C(O)NR'R", -SO<sub>2</sub>R', unsubstituted or substituted C<sub>1-8</sub> alkyl (optionally C4-8 [[1-8]] alkoxyalkyloxy, CH2CH2OCH2CH2OMe CH2CH2OMe) alkyl, unsubstituted or substituted C<sub>2-8</sub> alkenyl, unsubstituted or substituted C<sub>3-8</sub> alkynyl, unsubstituted or substituted C<sub>3-8</sub> cycloalkyl, unsubstituted or substituted 3- to 10-membered heterocyclyl.

- 28. (Previously presented) The modulator of claim 27, where  ${\sf R}^7$  is  ${\sf C}_{\sf 1-8}$  alkoxyalkyloxy.
- 29. (Original) The modulator of claim 27, where n is 1.
- 30. (Original) A modulator comprising one of the following formulae:

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19		69	

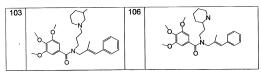
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22		72	\$0,50 0,50 0,50 0,50
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24		74	
25	->-10	75	
26	->-100	76	
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35	- Children	85	55,00
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101		104	
102	2	105	» ( S)



- 31. (Currently amended) A pharmaceutical composition comprising the modulator of claim 1 and a pharmaceutically acceptable carrier.
- 32. (Original) A pharmaceutical composition comprising the modulator of claim 27 and a pharmaceutically acceptable carrier.
- 33. (Withdrawn) A pharmaceutical composition comprising the modulator of claim 30 and a pharmaceutically acceptable carrier.
- 34. (Withdrawn) A pharmaceutical composition comprising a compound of the formulae:

Amendment dated January 6, 2008 Reply to Office Action of September 6, 2007 Appl. No. 10/743,281

and a pharmaceutically acceptable carrier.

- 35. (Withdrawn) A method of inhibiting the binding of chemokines I-TAC and/or SDF-1 to a CCXCKR2 receptor, comprising contacting the composition of claim 34 with a cell that expresses the CCXCKR2 receptor for a time sufficient to inhibit the binding of the chemokines to the CCXCKR2 receptor.
- 36. (Withdrawn) A method of inhibiting the binding of chemokines I-TAC and/or SDF-1 to a CCXCKR2 receptor, comprising contacting the modulator of claim 1 with a cell that expresses the CCXCKR2 receptor for a time sufficient to inhibit the binding of the chemokines to the CCXCKR2 receptor.
- 37. (Withdrawn) A method of treating cancer, comprising administering a therapeutically effective amount of the composition of claim 3234 34 to a cancer patient for a time sufficient to treat the cancer.
- 38. (Withdrawn) A method of treating cancer, comprising administering a therapeutically effective amount of the modulator of claim 1 to a cancer patient for a time sufficient to treat the cancer.